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=> d his

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O
L7         639687 S L6 OR NUCLEIC ACID#
L8              4 S L7 AND L5
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broad structure search

=> fil reg

FILE 'REGISTRY' ENTERED AT 10:17:42 ON 10 SEP 2001
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

STRUCTURE FILE UPDATES: 7 SEP 2001 HIGHEST RN 355367-45-2
DICTIONARY FILE UPDATES: 7 SEP 2001 HIGHEST RN 355367-45-2

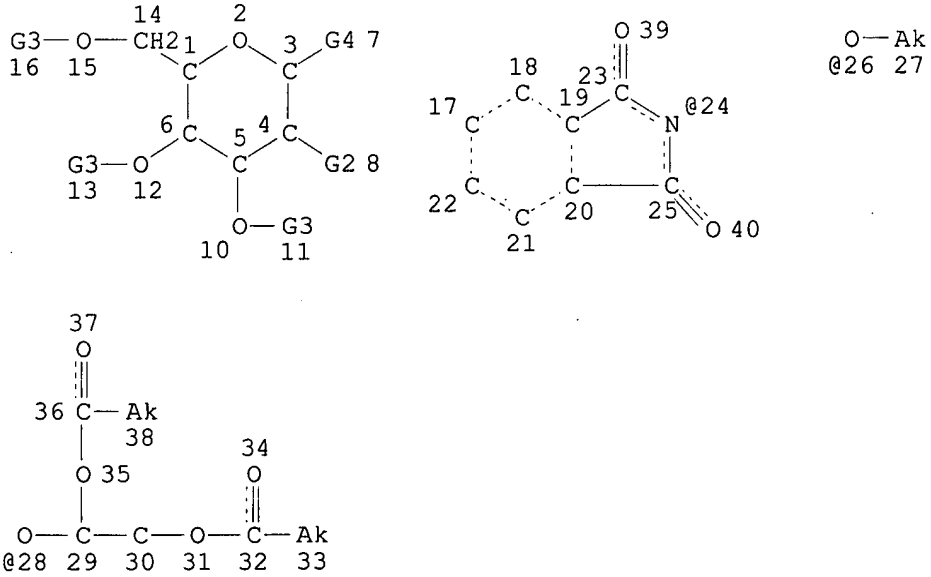
TSCA INFORMATION NOW CURRENT THROUGH January 11, 2001

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Structure search limits have been increased. See HELP SLIMIT for details.

=> d que stat l4

L3 STR



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VAR G3=H/C(O)CH3
VAR G4=BR/26/28
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CONNECT IS E1 RC AT 27
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CONNECT IS E1 RC AT 38
DEFAULT MLEVEL IS ATOM
DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:
RING(S) ARE ISOLATED OR EMBEDDED

NUMBER OF NODES IS 39

STEREO ATTRIBUTES: NONE
L4 149 SEA FILE=REGISTRY SSS FUL L3

100.0% PROCESSED 50974 ITERATIONS 149 ANSWERS
SEARCH TIME: 00.00.15

=> fil hcaplus

FILE 'HCAPLUS' ENTERED AT 10:17:52 ON 10 SEP 2001
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

FILE COVERS 1947 - 10 Sep 2001 VOL 135 ISS 12
FILE LAST UPDATED: 7 Sep 2001 (20010907/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

This file supports REG1stRY for direct browsing and searching of all substance data from the REGISTRY file. Enter HELP FIRST for more information.

HCAplus now provides online access to patents and literature covered in CA from 1947 to the present. On April 22, 2001, bibliographic information and abstracts were added for over 2.2 million references published in CA from 1947 to 1966.

'OBI' IS DEFAULT SEARCH FIELD FOR 'HCAPLUS' FILE

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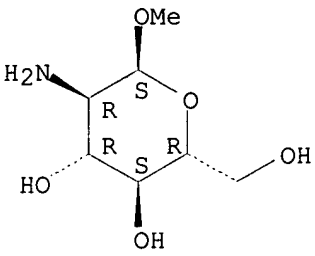
=> d his l5-

FILE 'HCAPLUS' ENTERED AT 10:13:40 ON 10 SEP 2001
L5 401 S L4
L6 611745 S NUCLEOTIDE# OR DNA OR RNA OR POLYNUCLEOTIDE# OR PLASMID# OR O
L7 639687 S L6 OR NUCLEIC ACID#
L8 4 S L7 AND L5

=> d .ca hitstr 1-4

L8 ANSWER 1 OF 4 HCAPLUS COPYRIGHT 2001 ACS
ACCESSION NUMBER: 2001:186432 HCAPLUS
DOCUMENT NUMBER: 134:340652
TITLE: A new strategy for the solid-phase synthesis of glycoconjugate biomolecules
AUTHOR(S): Di Fabio, Giovanni; De Capua, Antonia; De Napoli, Lorenzo; Montesarchio, Daniela; Piccialli, Gennaro; Rossi, Filomena; Benedetti, Ettore
CORPORATE SOURCE: Dipartimento di Chimica Organica e Biochimica, Universita degli Studi di Napoli "Federico II", Naples, 80126, Italy
SOURCE: Synlett (2001), (3), 341-344
CODEN: SYNLES; ISSN: 0936-5214
PUBLISHER: Georg Thieme Verlag
DOCUMENT TYPE: Journal
LANGUAGE: English
AB A simple and efficient bi-directional solid-phase synthesis, based on the use of a Tentagel solid support, functionalized with a suitably protected 2-amino sugar residue, is proposed for the prepn. of a variety of glycoconjugates, including glycopeptides and nucleo-glycopeptides.
CC 33-9 (Carbohydrates)
Section cross-reference(s): 34
ST **nucleotide** glycoconjugate glycopeptide Tentagel solid phase synthesis
IT Solid phase synthesis
(solid-phase synthesis of **nucleotide** glycopeptides)
IT Glycoconjugates
Glycopeptides
RL: SPN (Synthetic preparation); PREP (Preparation)
(solid-phase synthesis of **nucleotide** glycopeptides)
IT 108-30-5, Succinic anhydride, reactions **13374-29-3** 35661-60-0 98796-53-3 102212-98-6 313988-69-1
RL: RCT (Reactant)
(solid-phase synthesis of **nucleotide** glycopeptides)
IT 338801-55-1P 338801-56-2DP, Tentagel resin polymer support
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation)
(solid-phase synthesis of **nucleotide** glycopeptides)
IT 338801-57-3P 338801-58-4P 338801-59-5P 338801-60-8P
RL: SPN (Synthetic preparation); PREP (Preparation)
(solid-phase synthesis of **nucleotide** glycopeptides)
IT **13374-29-3**
RL: RCT (Reactant)
(solid-phase synthesis of **nucleotide** glycopeptides)
RN 13374-29-3 HCAPLUS
CN .alpha.-D-Glucopyranoside, methyl 2-amino-2-deoxy-, hydrochloride (7CI, 9CI) (CA INDEX NAME)

Absolute stereochemistry.



● HCl

REFERENCE COUNT: 23
REFERENCE(S):
(1) Adinolfi, M; Tetrahedron Lett 1996, V37, P5007 HCAPLUS
(2) Adinolfi, M; Tetrahedron Lett 1998, V39, P1953 HCAPLUS
(3) Adinolfi, M; Tetrahedron Lett 1999, V40, P2607 HCAPLUS
(4) Akhtar, S; Tetrahedron Lett 1995, V36, P7333 HCAPLUS
(7) de Kort, M; Eur J Org Chem 1999, P2337 HCAPLUS
ALL CITATIONS AVAILABLE IN THE RE FORMAT

L8 ANSWER 2 OF 4 HCAPLUS COPYRIGHT 2001 ACS
ACCESSION NUMBER: 1999:194158 HCAPLUS
DOCUMENT NUMBER: 130:242316
TITLE: Hydrophobic glycosylamine derivatives, compositions, and methods for their use
INVENTOR(S): Mumper, Russell J.; Tagliaferri, Frank
PATENT ASSIGNEE(S): Genemedicine, Inc., USA
SOURCE: PCT Int. Appl., 88 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 9912945	A2	19990318	WO 1998-US18888	19980908
WO 9912945	A3	19990819		
W: AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, EE, ES, FI, GB, GE, GH, GM, HR, HU, ID, IL, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, UA, UG, US, UZ, VN, YU, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
RW: GH, GM, KE, LS, MW, SD, SZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG				
AU 9893839	A1	19990329	AU 1998-93839	19980908
EP 1015465	A2	20000705	EP 1998-946932	19980908
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT,				

IE, FI
PRIORITY APPLN. INFO.: US 1997-58259 P 19970908
WO 1998-US18888 W 19980908
OTHER SOURCE(S): MARPAT 130:242316

AB The invention relates in part to hydrophobic glycosylamine derivs., methods for synthesizing hydrophobic derivs., compns. comprising these derivs., and methods for delivering macromols., such as proteins, peptides, lipids, carbohydrates, peptidomimetics, org. mols., and nucleic acids, to cells by administering these compns. The compds., compns., and methods of the invention are particularly useful for gene therapy and cancer treatment. Compns. contg. 1-mono-oleyl-.beta.-D-glucosamine or 1-monopalmityl-.beta.-D-glucosamine, plasmid DNA comprising an IL-2 gene, and DOPE reduced the growth rate in mice by 30% after 9 days and by 25% after 13 days, resp.

IC ICM C07H015-00

CC 63-6 (Pharmaceuticals)
Section cross-reference(s): 33

IT **Nucleotides**, biological studies
RL: THU (Therapeutic use); BIOL (Biological study); USES (Uses)
(analogs; compns. and synthesis of hydrophobic glycosylamine derivs. for delivery of macromol. compds. to cells)

IT Antitumor agents
Drug delivery systems
Gene therapy
Infusions (drug delivery systems)
Inhalants (drug delivery systems)
Injections (drug delivery systems)
Intravenous injections
Liposomes (drug delivery systems)
Oral drug delivery systems
Plasmids
Reducing agents
(compns. and synthesis of hydrophobic glycosylamine derivs. for delivery of macromol. compds. to cells)

IT **DNA**
Nucleic acids
Polynucleotides
RNA
RL: THU (Therapeutic use); BIOL (Biological study); USES (Uses)
(compns. and synthesis of hydrophobic glycosylamine derivs. for delivery of macromol. compds. to cells)

IT **221247-53-6P 221247-55-8P**
RL: PRP (Properties); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(compns. and synthesis of hydrophobic glycosylamine derivs. for delivery of macromol. compds. to cells)

IT 143-28-2, Oleyl alcohol 36653-82-4, Hexadecanol **63000-69-1**
138395-62-7
RL: RCT (Reactant)
(compns. and synthesis of hydrophobic glycosylamine derivs. for delivery of macromol. compds. to cells)

IT **221247-52-5P 221247-54-7P**
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation)
(compns. and synthesis of hydrophobic glycosylamine derivs. for delivery of macromol. compds. to cells)

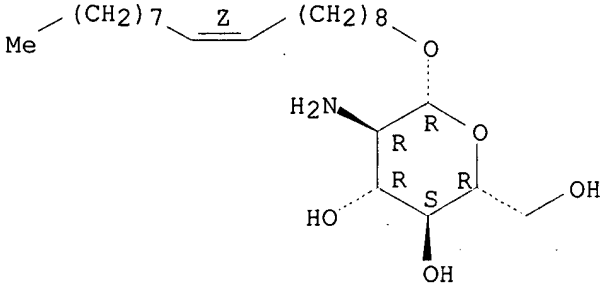
IT **221247-53-6P 221247-55-8P**
RL: PRP (Properties); SPN (Synthetic preparation); THU (Therapeutic use);

BIOL (Biological study); PREP (Preparation); USES (Uses)
(compns. and synthesis of hydrophobic glycosylamine derivs. for delivery of macromol. compds. to cells)

RN 221247-53-6 HCAPLUS

CN .beta.-D-Glucopyranoside, (9Z)-9-octadecenyl 2-amino-2-deoxy- (9CI) (CA INDEX NAME)

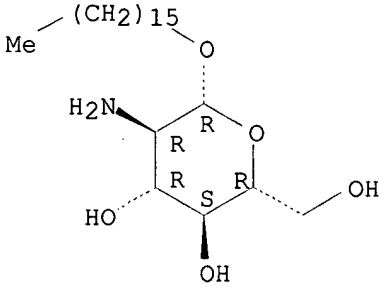
Absolute stereochemistry.
Double bond geometry as shown.



RN 221247-55-8 HCAPLUS

CN .beta.-D-Glucopyranoside, hexadecyl 2-amino-2-deoxy- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

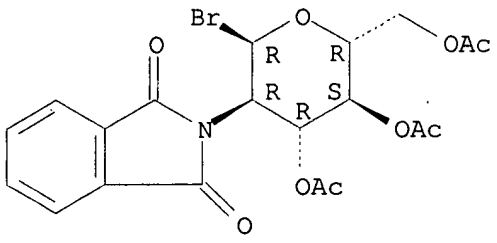


IT **63000-69-1 138395-62-7**
RL: RCT (Reactant)
(compns. and synthesis of hydrophobic glycosylamine derivs. for delivery of macromol. compds. to cells)

RN 63000-69-1 HCAPLUS

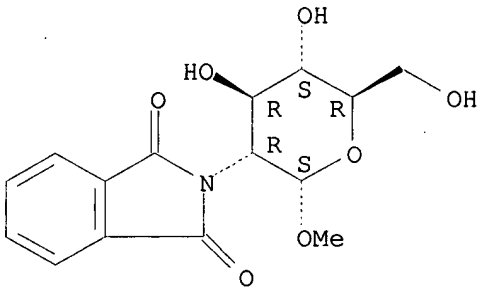
CN .alpha.-D-Glucopyranosyl bromide, 2-deoxy-2-(1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)-, 3,4,6-triacetate (9CI) (CA INDEX NAME)

Absolute stereochemistry.



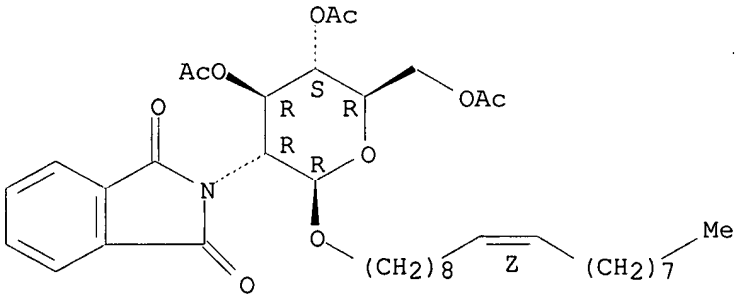
RN 138395-62-7 HCAPLUS
CN .alpha.-D-Glucopyranoside, methyl 2-deoxy-2-(1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



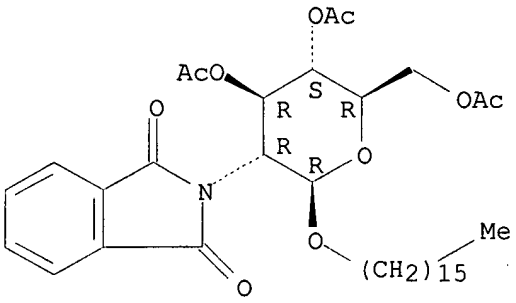
IT **221247-52-5P 221247-54-7P**
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation)
(compns. and synthesis of hydrophobic glycosylamine derivs. for delivery of macromol. compds. to cells)
RN 221247-52-5 HCAPLUS
CN .beta.-D-Glucopyranoside, (9Z)-9-octadecenyl 2-deoxy-2-(1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)-, 3,4,6-triacetate (9CI) (CA INDEX NAME)

Absolute stereochemistry.
Double bond geometry as shown.



RN 221247-54-7 HCAPLUS
CN .beta.-D-Glucopyranoside, hexadecyl 2-deoxy-2-(1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)-, 3,4,6-triacetate (9CI) (CA INDEX NAME)

Absolute stereochemistry.



L8 ANSWER 3 OF 4 HCAPLUS COPYRIGHT 2001 ACS
ACCESSION NUMBER: 1995:441303 HCAPLUS
DOCUMENT NUMBER: 122:205195
TITLE: Bivalent sialyl Lewis X (SLex) saccharides to inhibit selectin-mediated cell adhesion
INVENTOR(S): Gaeta, Federico C. A.; DeFrees, Shawn A.
PATENT ASSIGNEE(S): Cytel Corp., USA
SOURCE: PCT Int. Appl., 105 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 2
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 9503059	A1	19950202	WO 1994-US8384	19940721
W: AU, BG, CA, CN, CZ, FI, HU, JP, KR, MN, NO, NZ, PL, RO, RU, SK, UA, VN				
RW: AT, BE, CH, DE, DK, ES, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, ML, MR, NE, SN, TD, TG				
US 5559103	A	19960924	US 1994-278020	19940720
AU 9474046	A1	19950220	AU 1994-74046	19940721
JP 09500683	T2	19970121	JP 1994-505381	19940721

PRIORITY APPLN. INFO.: US 1993-95657 19930721
US 1994-278020 19940720
WO 1994-US8384 19940721

OTHER SOURCE(S): MARPAT 122:205195

AB Bivalent SLex saccharide derivs. are provided that inhibit binding of cells expressing a surface selectin receptor (e.g. blood platelets and vascular endothelial cells) to cells which express SLex on their surfaces (e.g. leukocytes). Pharmaceutical compns. comprising these saccharides, processes for making and using them, and methods for synthesis of the saccharides are also disclosed.

IC ICM A61K031-715
CC 1-8 (Pharmacology)
Section cross-reference(s): 33, 63
IT **Plasmid** and Episome
(pCDNAI-sol-E-selectin; bivalent sialyl Lewis X (SLex) saccharides to inhibit selectin-mediated cell adhesion)
IT 98-88-4, Benzoyl chloride 3019-71-4, Trichloroacetyl isocyanate

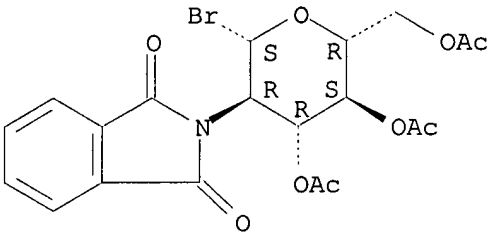
10028-45-2 28053-08-9 28605-65-4 101833-22-1 148296-47-3
RL: RCT (Reactant)
(bivalent sialyl Lewis X (SLex) saccharides to inhibit selectin-mediated cell adhesion)

IT **10028-45-2**
RL: RCT (Reactant)
(bivalent sialyl Lewis X (SLex) saccharides to inhibit selectin-mediated cell adhesion)

RN 10028-45-2 HCAPLUS

CN .beta.-D-Glucopyranosyl bromide, 2-deoxy-2-(1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)-, 3,4,6-triacetate (9CI) (CA INDEX NAME)

Absolute stereochemistry.



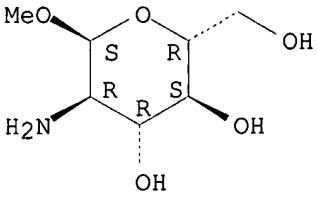
L8 ANSWER 4 OF 4 HCAPLUS COPYRIGHT 2001 ACS
ACCESSION NUMBER: 1994:548341 HCAPLUS
DOCUMENT NUMBER: 121:148341
TITLE: Structure-activity relationships in the induction of single-strand breakage in **plasmid** pBR322 **DNA** by amino sugars and derivatives
AUTHOR(S): Kashige, Nobuhiro; Yamaguchi, Tadatoshi; Ohtakara, Akira; Mitsutomi, Masaru; Brimacombe, John S.; Miake, Fumio; Watanabe, Kenji
CORPORATE SOURCE: Faculty of Pharmaceutical Sciences, Fukuoka University, Nanakuma, Jonan-ku, Fukuoka, 814-01, Japan
SOURCE: Carbohydr. Res. (1994), 257(2), 285-91
CODEN: CRBRAT; ISSN: 0008-6215
DOCUMENT TYPE: Journal
LANGUAGE: English

AB Structure-activity relationships in the induction of strand breakage in plasmid pBR322 DNA by amino sugars and their derivs. were investigated using agarose gel electrophoresis. The coexistence of a potential free aldehyde group at the C-1 position and a free amino group at the C-2 position in the mols. was indispensable for the display of DNA strand-breaking activity in both mono- and oligo-aminosaccharides. The activity was increased by the introduction of an acidic group, esp. a phosphate group, at the C-6 position. The activity was also increased by the addn. of Cu2+. The order of activity of the amino monosaccharides tested was D-isoglucosamine > D-mannosamine > D-galactosamine > D-glucosamine, and it is suggested that this order is correlated with the portion of acyclic (aldehydo) form in the soln. of each sugar. The possible chem. basis for DNA strand breakage by amino sugars is discussed.

CC 1-3 (Pharmacology)
Section cross-reference(s): 4

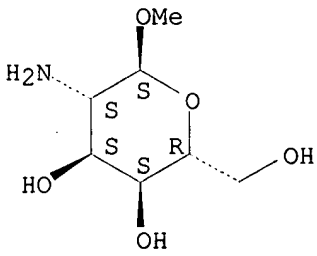
ST amino sugar **DNA** strand breakage structure
IT Molecular structure-biological activity relationship
(**DNA**-cleaving, of amino sugars and derivs.)
IT 50-99-7, D-Glucose, biological studies 154-17-6, 2-Deoxy-D-glucose
298-08-8, .alpha.-Aminoacetone 499-14-9, Chondrosine 576-44-3,
D-Kanosamine 577-76-4, Chitobiose 1811-31-0, N-Acetyl-D-galactosamine
3416-24-8, D-Glucosamine 3615-17-6, N-Acetyl-D-mannosamine 3616-42-0,
D-Glucosamine 6-phosphate 3646-68-2, D-Glucosaminic acid 4429-04-3,
D-Isoglucosamine 4607-22-1 **4704-14-7**, Methyl
2-amino-2-deoxy-.alpha.-D-glucopyranoside 5155-47-5, Methyl
6-amino-6-deoxy-.alpha.-D-glucopyranoside 5567-52-2, Chitotetraose
7512-17-6, N-Acetyl-D-glucosamine 7535-00-4, D-Galactosamine
14307-02-9, D-Mannosamine 14635-95-1 19889-76-0 32385-07-2
35812-81-8 40879-84-3 41708-93-4, Chitotriose 41708-94-5,
Chitopentaose 41708-95-6, Chitohexaose 50692-69-8, Benzyl
2-amino-2-deoxy-.alpha.-D-glucopyranoside 53574-53-1 **63122-16-7**
66954-08-3 76330-20-6 91674-26-9 138430-53-2 156304-79-9
156304-80-2 156304-81-3
RL: BIOL (Biological study)
(**DNA** single-strand breakage by, structure in relation to)
IT **4704-14-7**, Methyl 2-amino-2-deoxy-.alpha.-D-glucopyranoside
63122-16-7
RL: BIOL (Biological study)
(**DNA** single-strand breakage by, structure in relation to)
RN 4704-14-7 HCAPLUS
CN .alpha.-D-Glucopyranoside, methyl 2-amino-2-deoxy- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



RN 63122-16-7 HCAPLUS
CN .alpha.-D-Altropyranoside, methyl 2-amino-2-deoxy- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



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L4             24 SEA FILE=REGISTRY SUB=L2 SSS FUL L3
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L5         501021 S NC4-C6/ES
L6             14 S L5 AND L4
L7              5 S L4 AND C>23
L8             18 S L4 AND BR/ELS
L9             12 S L8 AND L6

     FILE 'HCAPLUS' ENTERED AT 10:08:32 ON 10 SEP 2001
L10             6 S L7
L11           194 S L9
L12             1 S L10 AND L11
L13             5 S L10 NOT L12
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narrowed structure search

```
=> fil reg

FILE 'REGISTRY' ENTERED AT 10:09:54 ON 10 SEP 2001
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
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STRUCTURE FILE UPDATES:    7 SEP 2001  HIGHEST RN 355367-45-2
DICTIONARY FILE UPDATES:   7 SEP 2001  HIGHEST RN 355367-45-2

TSCA INFORMATION NOW CURRENT THROUGH January 11, 2001

  Please note that search-term pricing does apply when
  conducting SmartSELECT searches.

Structure search limits have been increased.  See HELP SLIMIT
for details.

=> d his l1-19


     (FILE 'HCAPLUS' ENTERED AT 10:01:46 ON 10 SEP 2001)
               DEL HIS Y

     FILE 'REGISTRY' ENTERED AT 10:03:17 ON 10 SEP 2001
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L1                STR
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L3                STR
L4             24 SEA FILE=REGISTRY SUB=L2 SSS FUL L3
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L5         501021 S NC4-C6/ES
L6             14 S L5 AND L4
L7              5 S L4 AND C>23
L8             18 S L4 AND BR/ELS
L9             12 S L8 AND L6

=> d que stat l4

L1                STR
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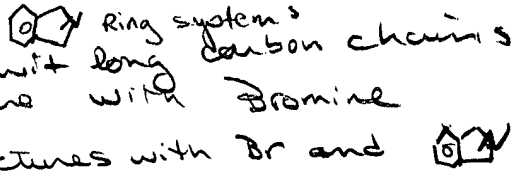
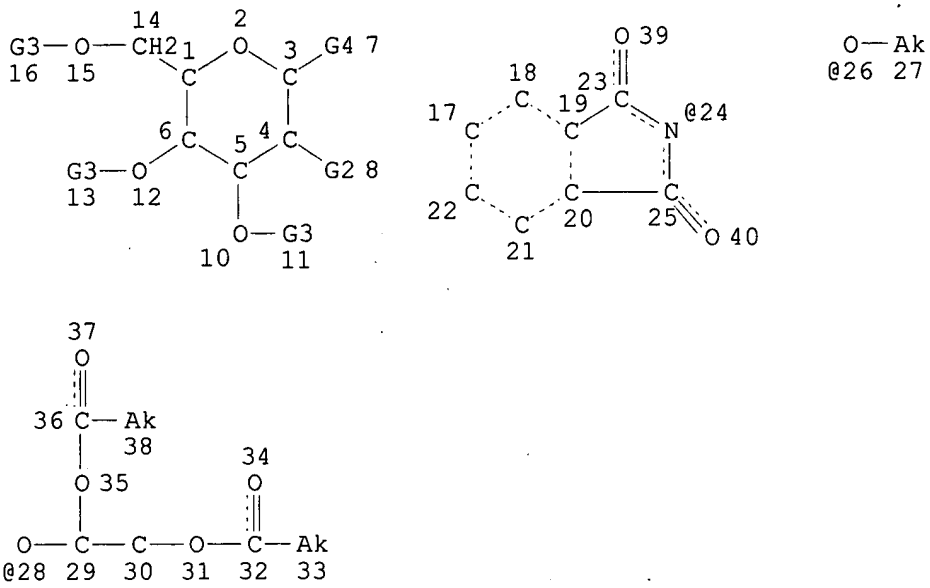
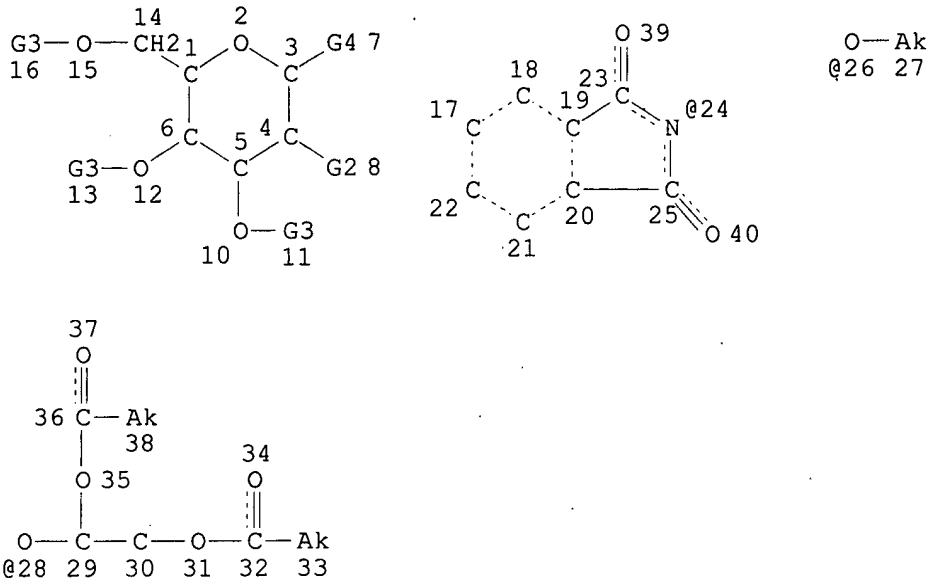




VAR G2=NH2/24
VAR G3=H/C(O)CH3
VAR G4=BR/26/28
NODE ATTRIBUTES:
CONNECT IS E1 RC AT 27
CONNECT IS E1 RC AT 33
CONNECT IS E1 RC AT 38
DEFAULT MLEVEL IS ATOM
DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:
RING(S) ARE ISOLATED OR EMBEDDED
NUMBER OF NODES IS 39

STEREO ATTRIBUTES: NONE
L2 (149)SEA FILE=REGISTRY SSS FUL L1
L3 STR



VAR G2=NH2/24
VAR G3=H/C(O)CH3
VAR G4=BR/26/28
NODE ATTRIBUTES:
CONNECT IS E1 RC AT 27
CONNECT IS E1 RC AT 33
CONNECT IS E1 RC AT 38
DEFAULT MLEVEL IS ATOM
DEFAULT ECLEVEL IS LIMITED
ECOUNT IS M16-X18 C AT 27
ECOUNT IS M16-X18 C AT 33
ECOUNT IS M16-X18 C AT 38

GRAPH ATTRIBUTES:
RING(S) ARE ISOLATED OR EMBEDDED
NUMBER OF NODES IS 39

STEREO ATTRIBUTES: NONE
L4 24 SEA FILE=REGISTRY SUB=L2 SSS FUL L3

100.0% PROCESSED 149 ITERATIONS 24 ANSWERS
SEARCH TIME: 00.00.02

=> d his l4-l9

(FILE 'REGISTRY' ENTERED AT 10:03:17 ON 10 SEP 2001)
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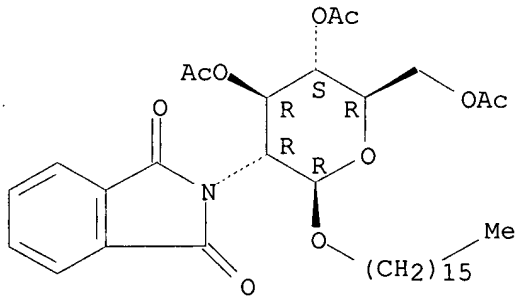
L5 501021 S NC4-C6/ES
L6 14 S L5 AND L4

L7 5 S L4 AND C>23
L8 18 S L4 AND BR/ELS
L9 12 S L8 AND L6

=> d ide can 17 1-5;d ide can 19 1-12

L7 ANSWER 1 OF 5 REGISTRY COPYRIGHT 2001 ACS
RN 221247-54-7 REGISTRY
CN .beta.-D-Glucopyranoside, hexadecyl 2-deoxy-2-(1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)-, 3,4,6-triacetate (9CI) (CA INDEX NAME)
FS STEREOSEARCH
MF **C36 H53 N O10**
SR CA
LC STN Files: CA, CAPLUS, TOXLIT

Absolute stereochemistry.

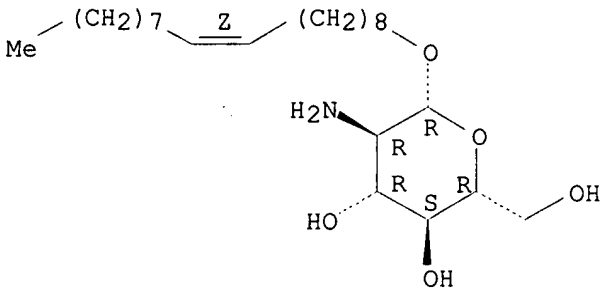


1 REFERENCES IN FILE CA (1967 TO DATE)
1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 130:242316

L7 ANSWER 2 OF 5 REGISTRY COPYRIGHT 2001 ACS
RN 221247-53-6 REGISTRY
CN .beta.-D-Glucopyranoside, (9Z)-9-octadecenyl 2-amino-2-deoxy- (9CI) (CA INDEX NAME)
FS STEREOSEARCH
MF **C24 H47 N O5**
SR CA
LC STN Files: CA, CAPLUS, TOXLIT

Absolute stereochemistry.
Double bond geometry as shown.

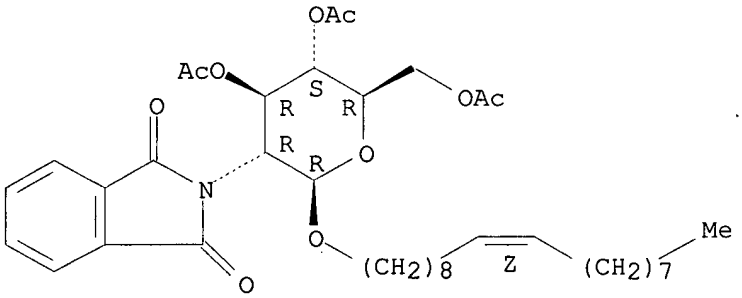


1 REFERENCES IN FILE CA (1967 TO DATE)
1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 130:242316

L7 ANSWER 3 OF 5 REGISTRY COPYRIGHT 2001 ACS
RN 221247-52-5 REGISTRY
CN .beta.-D-Glucopyranoside, (9Z)-9-octadecenyl 2-deoxy-2-(1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)-, 3,4,6-triacetate (9CI) (CA INDEX NAME)
FS STEREOSEARCH
MF **C38 H55 N O10**
SR CA
LC STN Files: CA, CAPLUS, TOXLIT

Absolute stereochemistry.
Double bond geometry as shown.

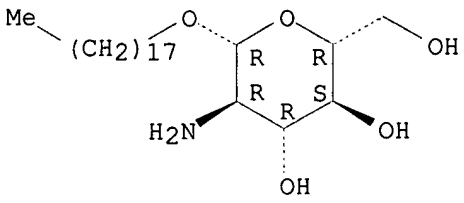


1 REFERENCES IN FILE CA (1967 TO DATE)
1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 130:242316

L7 ANSWER 4 OF 5 REGISTRY COPYRIGHT 2001 ACS
RN 159405-31-9 REGISTRY
CN .beta.-D-Glucopyranoside, octadecyl 2-amino-2-deoxy- (9CI) (CA INDEX NAME)
FS STEREOSEARCH
MF **C24 H49 N O5**
SR CA
LC STN Files: CA, CAPLUS

Absolute stereochemistry.



4 REFERENCES IN FILE CA (1967 TO DATE)
4 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 131:333590

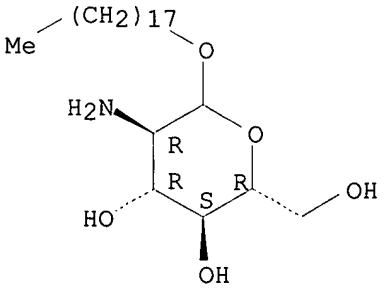
REFERENCE 2: 128:305665

REFERENCE 3: 124:176728

REFERENCE 4: 121:312595

L7 ANSWER 5 OF 5 REGISTRY COPYRIGHT 2001 ACS
RN 134227-32-0 REGISTRY
CN D-Glucopyranoside, octadecyl 2-amino-2-deoxy-, hydrochloride (9CI) (CA INDEX NAME)
FS STEREOSEARCH
MF **C24 H49 N O5 . Cl H**
SR CA
LC STN Files: CA, CAPLUS, TOXLIT

Absolute stereochemistry.



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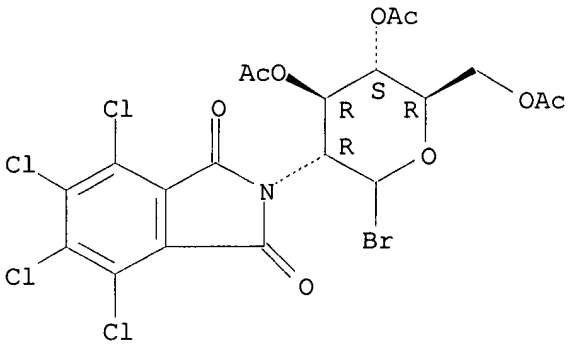
1 REFERENCES IN FILE CA (1967 TO DATE)
1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 115:15307

L9 ANSWER 1 OF 12 REGISTRY COPYRIGHT 2001 ACS
RN 189218-64-2 REGISTRY
CN D-Glucopyranosyl bromide, 2-deoxy-2-(4,5,6,7-tetrachloro-1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)-, 3,4,6-triacetate (9CI) (CA INDEX NAME)
FS STEREOSEARCH
MF **C20 H16 Br Cl4 N O9**
SR CA
LC STN Files: CA, CAPLUS, CASREACT

Absolute stereochemistry.



3 REFERENCES IN FILE CA (1967 TO DATE)
3 REFERENCES IN FILE CAPLUS (1967 TO DATE)

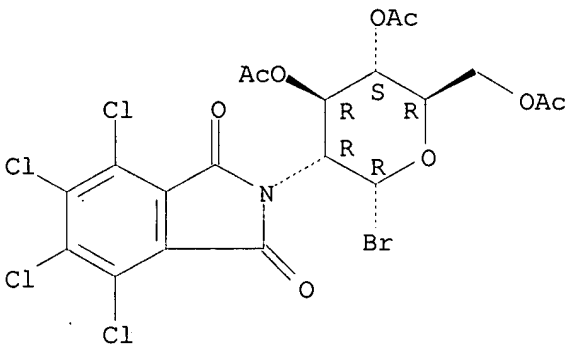
REFERENCE 1: 134:5080

REFERENCE 2: 129:331058

REFERENCE 3: 126:305712

L9 ANSWER 2 OF 12 REGISTRY COPYRIGHT 2001 ACS
RN 180778-39-6 REGISTRY
CN .alpha.-D-Glucopyranosyl bromide, 2-deoxy-2-(4,5,6,7-tetrachloro-1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)-, 3,4,6-triacetate (9CI) (CA INDEX NAME)
FS STEREOSEARCH
MF **C20 H16 Br Cl4 N O9**
SR CA
LC STN Files: CA, CAPLUS, CASREACT

Absolute stereochemistry.

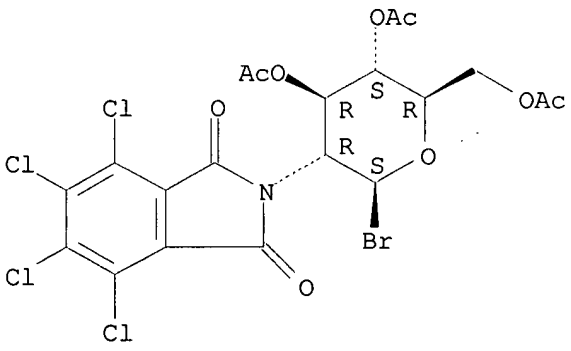


1 REFERENCES IN FILE CA (1967 TO DATE)
1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 125:196181

L9 ANSWER 3 OF 12 REGISTRY COPYRIGHT 2001 ACS
RN 180778-38-5 REGISTRY
CN .beta.-D-Glucopyranosyl bromide, 2-deoxy-2-(4,5,6,7-tetrachloro-1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)-, 3,4,6-triacetate (9CI) (CA INDEX NAME)
FS STEREOSEARCH
MF **C20 H16 Br Cl4 N O9**
SR CA
LC STN Files: CA, CAPLUS, CASREACT

Absolute stereochemistry.



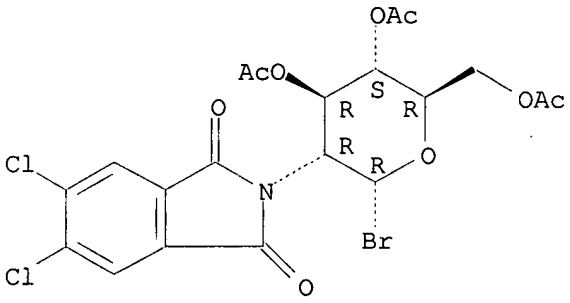
1 REFERENCES IN FILE CA (1967 TO DATE)
1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 125:196181

L9 ANSWER 4 OF 12 REGISTRY COPYRIGHT 2001 ACS
RN 175229-72-8 REGISTRY
CN .alpha.-D-Glucopyranosyl bromide, 2-deoxy-2-(5,6-dichloro-1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)-, 3,4,6-triacetate (9CI) (CA INDEX NAME)
FS STEREOSEARCH

MF **C20 H18 Br Cl2 N O9**
SR CA
LC STN Files: CA, CAPLUS

Absolute stereochemistry.

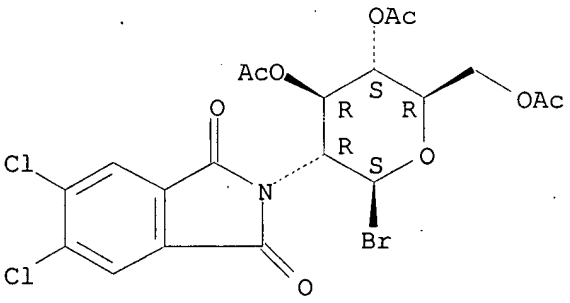


1 REFERENCES IN FILE CA (1967 TO DATE)
1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 124:261532

L9 ANSWER 5 OF 12 REGISTRY COPYRIGHT 2001 ACS
RN 175229-71-7 REGISTRY
CN .beta.-D-Glucopyranosyl bromide, 2-deoxy-2-(5,6-dichloro-1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)-, 3,4,6-triacetate (9CI) (CA INDEX NAME)
FS STEREOSEARCH
MF **C20 H18 Br Cl2 N O9**
SR CA
LC STN Files: CA, CAPLUS

Absolute stereochemistry. Rotation (+).



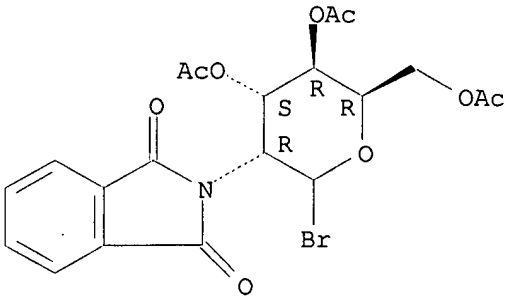
1 REFERENCES IN FILE CA (1967 TO DATE)
1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 124:261532

L9 ANSWER 6 OF 12 REGISTRY COPYRIGHT 2001 ACS
RN 171234-13-2 REGISTRY
CN D-Gulopyranosyl bromide, 2-deoxy-2-(1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)-, 3,4,6-triacetate (9CI) (CA INDEX NAME)

FS STEREOSEARCH
MF **C20 H20 Br N O9**
SR CA
LC STN Files: CA, CAPLUS

Absolute stereochemistry.



3 REFERENCES IN FILE CA (1967 TO DATE)
3 REFERENCES IN FILE CAPLUS (1967 TO DATE)

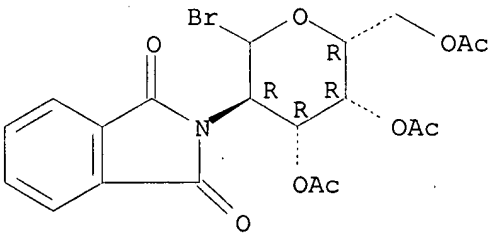
REFERENCE 1: 125:222299

REFERENCE 2: 124:56497

REFERENCE 3: 124:9205

L9 ANSWER 7 OF 12 REGISTRY COPYRIGHT 2001 ACS
RN 110455-06-6 REGISTRY
CN D-Galactopyranosyl bromide,
2-deoxy-2-(1,3-dihydro-1,3-dioxo-2H-isoindol-2-
yl)-, 3,4,6-triacetate (9CI) (CA INDEX NAME)
FS STEREOSEARCH
MF **C20 H20 Br N O9**
SR CA
LC STN Files: BEILSTEIN*, CA, CAPLUS, CASREACT, CHEMINFORMRX, TOXLIT,
USPATFULL
(*File contains numerically searchable property data)

Absolute stereochemistry.



8 REFERENCES IN FILE CA (1967 TO DATE)
8 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 122:133606

REFERENCE 2: 114:185935

REFERENCE 3: 113:59707

REFERENCE 4: 113:41172

REFERENCE 5: 111:23841

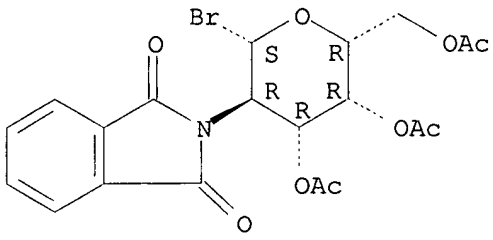
REFERENCE 6: 110:75984

REFERENCE 7: 109:223663

REFERENCE 8: 107:154687

L9 ANSWER 8 OF 12 REGISTRY COPYRIGHT 2001 ACS
RN 90458-06-3 REGISTRY
CN .beta.-D-Galactopyranosyl bromide, 2-deoxy-2-(1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)-, 3,4,6-triacetate (9CI) (CA INDEX NAME)
FS STEREOSEARCH
MF **C20 H20 Br N O9**
LC STN Files: BEILSTEIN*, CA, CAPLUS, CASREACT, CHEMINFORMRX, USPATFULL
(*File contains numerically searchable property data)

Absolute stereochemistry.



13 REFERENCES IN FILE CA (1967 TO DATE)
13 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 124:202848

REFERENCE 2: 123:33556

REFERENCE 3: 115:9257

REFERENCE 4: 113:172491

REFERENCE 5: 109:129528

REFERENCE 6: 108:6300

REFERENCE 7: 107:176347

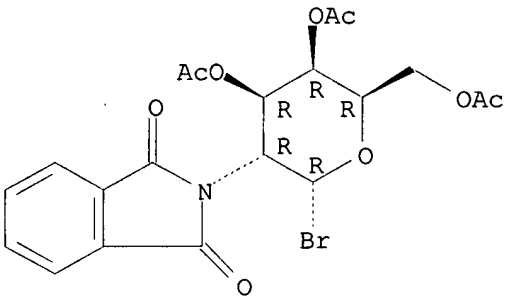
REFERENCE 8: 107:134594

REFERENCE 9: 106:120180

REFERENCE 10: 106:117683

L9 ANSWER 9 OF 12 REGISTRY COPYRIGHT 2001 ACS
RN 81704-03-2 REGISTRY
CN .alpha.-D-Galactopyranosyl bromide, 2-deoxy-2-(1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)-, 3,4,6-triacetate (9CI) (CA INDEX NAME)
FS STEREOSEARCH
MF **C20 H20 Br N O9**
LC STN Files: BEILSTEIN*, CA, CAPLUS, CASREACT, CHEMINFORMRX, TOXLIT, USPATFULL
(*File contains numerically searchable property data)

Absolute stereochemistry.



20 REFERENCES IN FILE CA (1967 TO DATE)
20 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 131:272136

REFERENCE 2: 131:157896

REFERENCE 3: 130:139551

REFERENCE 4: 127:5280

REFERENCE 5: 126:157708

REFERENCE 6: 121:158031

REFERENCE 7: 120:245690

REFERENCE 8: 120:54898

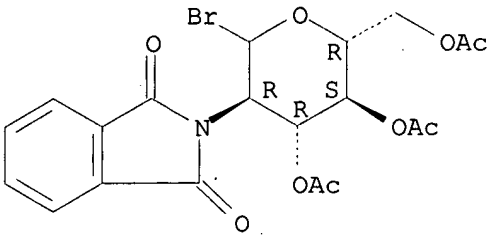
REFERENCE 9: 119:265156

REFERENCE 10: 117:111984

L9 ANSWER 10 OF 12 REGISTRY COPYRIGHT 2001 ACS
RN 70831-94-6 REGISTRY
CN D-Glucopyranosyl bromide, 2-deoxy-2-(1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)-, 3,4,6-triacetate (9CI) (CA INDEX NAME)
FS STEREOSEARCH
MF **C20 H20 Br N O9**
LC STN Files: BEILSTEIN*, CA, CAPLUS, CASREACT, CHEMINFORMRX, TOXLIT,

USPATFULL
(*File contains numerically searchable property data)

Absolute stereochemistry.



42 REFERENCES IN FILE CA (1967 TO DATE)
42 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 133:74212

REFERENCE 2: 131:45016

REFERENCE 3: 131:19212

REFERENCE 4: 129:316471

REFERENCE 5: 127:319184

REFERENCE 6: 126:168260

REFERENCE 7: 126:157708

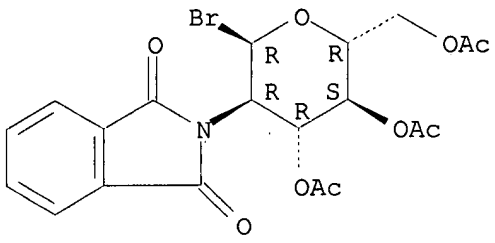
REFERENCE 8: 125:329206

REFERENCE 9: 125:185857

REFERENCE 10: 123:199268

L9 ANSWER 11 OF 12 REGISTRY COPYRIGHT 2001 ACS
RN 63000-69-1 REGISTRY
CN .alpha.-D-Glucopyranosyl bromide, 2-deoxy-2-(1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)-, 3,4,6-triacetate (9CI) (CA INDEX NAME)
FS STEREOSEARCH
MF **C20 H20 Br N O9**
LC STN Files: BEILSTEIN*, CA, CAPLUS, CASREACT, CHEMINFORMRX, TOXLIT, USPATFULL
(*File contains numerically searchable property data)

Absolute stereochemistry.



24 REFERENCES IN FILE CA (1967 TO DATE)
24 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 133:362913

REFERENCE 2: 130:242316

REFERENCE 3: 129:343650

REFERENCE 4: 126:47484

REFERENCE 5: 126:8452

REFERENCE 6: 125:301540

REFERENCE 7: 125:196166

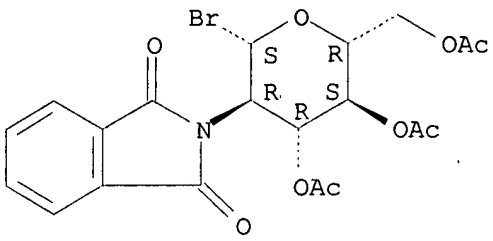
REFERENCE 8: 125:87030

REFERENCE 9: 123:33556

REFERENCE 10: 122:106316

L9 ANSWER 12 OF 12 REGISTRY COPYRIGHT 2001 ACS
RN 10028-45-2 REGISTRY
CN .beta.-D-Glucopyranosyl bromide, 2-deoxy-2-(1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)-, 3,4,6-triacetate (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Glucopyranosyl bromide, 2-deoxy-2-phthalimido-, 3,4,6-triacetate, .beta.-D- (6CI, 7CI, 8CI)
OTHER NAMES:
CN 2-Deoxy-2-phthalimido-3,4,6-tri-O-acetyl-.beta.-D-glucopyranosyl bromide
FS STEREOSEARCH
MF **C20 H20 Br N O9**
LC STN Files: BEILSTEIN*, CA, CAOLD, CAPLUS, CASREACT, CHEMCATS, CHEMINFORMRX, TOXLIT, USPATFULL
(*File contains numerically searchable property data)

Absolute stereochemistry.



103 REFERENCES IN FILE CA (1967 TO DATE)
103 REFERENCES IN FILE CAPLUS (1967 TO DATE)
4 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

REFERENCE 1: 134:42343

REFERENCE 2: 132:222783

REFERENCE 3: 131:72771

REFERENCE 4: 129:343648

REFERENCE 5: 129:290338

REFERENCE 6: 128:48415

REFERENCE 7: 126:238722

REFERENCE 8: 126:47484

REFERENCE 9: 125:329281

REFERENCE 10: 125:329158

=> fil hcaplus

FILE 'HCAPLUS' ENTERED AT 10:11:01 ON 10 SEP 2001
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

FILE COVERS 1947 - 10 Sep 2001 VOL 135 ISS 12
FILE LAST UPDATED: 7 Sep 2001 (20010907/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

This file supports REG1stRY for direct browsing and searching of all substance data from the REGISTRY file. Enter HELP FIRST for

more information.

HCAplus now provides online access to patents and literature covered in CA from 1947 to the present. On April 22, 2001, bibliographic information and abstracts were added for over 2.2 million references published in CA from 1947 to 1966.

'OBI' IS DEFAULT SEARCH FIELD FOR 'HCAPLUS' FILE

=> s hid l10-

L10- RANGE NOT VALID

=> d his l10-

(FILE 'HCAPLUS' ENTERED AT 10:08:32 ON 10 SEP 2001)
L10 6 S L7
L11 194 S L9
L12 1 S L10 AND L11
L13 5 S L10 NOT L12

FILE 'REGISTRY' ENTERED AT 10:09:54 ON 10 SEP 2001

FILE 'HCAPLUS' ENTERED AT 10:11:01 ON 10 SEP 2001

=> d .ca hitstr l12;d .ca hitstr l13 1-5

L12 ANSWER 1 OF 1 HCAPLUS COPYRIGHT 2001 ACS
ACCESSION NUMBER: 1999:194158 HCAPLUS
DOCUMENT NUMBER: 130:242316
TITLE: Hydrophobic glycosylamine derivatives, compositions, and methods for their use
INVENTOR(S): Mumper, Russell J.; Tagliaferri, Frank
PATENT ASSIGNEE(S): Genemedicine, Inc., USA
SOURCE: PCT Int. Appl., 88 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 9912945	A2	19990318	WO 1998-US18888	19980908
WO 9912945	A3	19990819		

W: AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, EE, ES, FI, GB, GE, GH, GM, HR, HU, ID, IL, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, UA, UG, US, UZ, VN, YU, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM
RW: GH, GM, KE, LS, MW, SD, SZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG

AU 9893839	A1	19990329	AU 1998-93839	19980908
EP 1015465	A2	20000705	EP 1998-946932	19980908

R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, FI

PRIORITY APPLN. INFO.: US 1997-58259 P 19970908
WO 1998-US18888 W 19980908

OTHER SOURCE(S): MARPAT 130:242316

AB The invention relates in part to hydrophobic glycosylamine derivs., methods for synthesizing hydrophobic derivs., compns. comprising these derivs., and methods for delivering macromols., such as proteins, peptides, lipids, carbohydrates, peptidomimetics, org. mols., and nucleic acids, to cells by administering these compns. The compds., compns., and methods of the invention are particularly useful for gene therapy and cancer treatment. Compns. contg. 1-mono-oleyl-.beta.-D-glucosamine or 1-monopalmityl-.beta.-D-glucosamine, plasmid DNA comprising an IL-2 gene, and DOPE reduced the growth rate in mice by 30% after 9 days and by 25% after 13 days, resp.

IC ICM C07H015-00

CC 63-6 (Pharmaceuticals)
Section cross-reference(s): 33

IT **221247-53-6P** 221247-55-8P
RL: PRP (Properties); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(compns. and synthesis of hydrophobic glycosylamine derivs. for delivery of macromol. compds. to cells)

IT 143-28-2, Oleyl alcohol 36653-82-4, Hexadecanol **63000-69-1** 138395-62-7
RL: RCT (Reactant)
(compns. and synthesis of hydrophobic glycosylamine derivs. for delivery of macromol. compds. to cells)

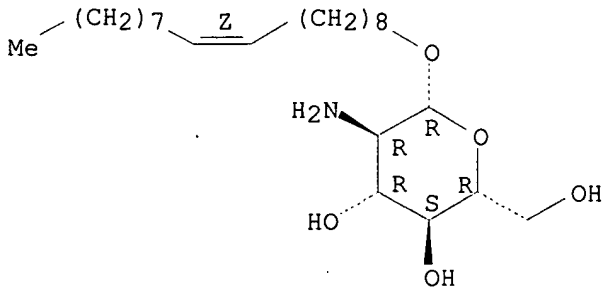
IT **221247-52-5P 221247-54-7P**
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation)
(compns. and synthesis of hydrophobic glycosylamine derivs. for delivery of macromol. compds. to cells)

IT **221247-53-6P**
RL: PRP (Properties); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(compns. and synthesis of hydrophobic glycosylamine derivs. for delivery of macromol. compds. to cells)

RN 221247-53-6 HCAPLUS

CN .beta.-D-Glucopyranoside, (9Z)-9-octadecenyl 2-amino-2-deoxy- (9CI) (CA INDEX NAME)

Absolute stereochemistry.
Double bond geometry as shown.

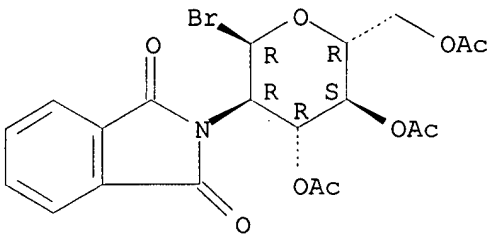


IT **63000-69-1**

RL: RCT (Reactant)
(compns. and synthesis of hydrophobic glycosylamine derivs. for delivery of macromol. compds. to cells)

RN 63000-69-1 HCAPLUS
CN .alpha.-D-Glucopyranosyl bromide, 2-deoxy-2-(1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)-, 3,4,6-triacetate (9CI) (CA INDEX NAME)

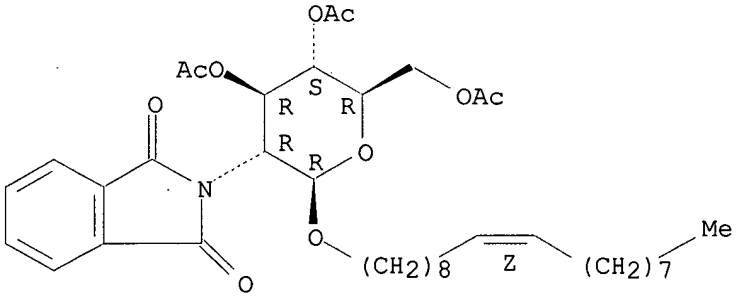
Absolute stereochemistry.



IT **221247-52-5P 221247-54-7P**
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation)
(compns. and synthesis of hydrophobic glycosylamine derivs. for delivery of macromol. compds. to cells)

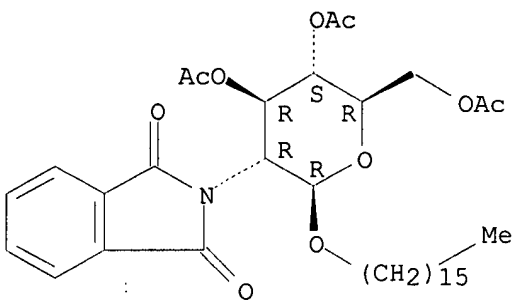
RN 221247-52-5 HCAPLUS
CN .beta.-D-Glucopyranoside, (9Z)-9-octadecenyl 2-deoxy-2-(1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)-, 3,4,6-triacetate (9CI) (CA INDEX NAME)

Absolute stereochemistry.
Double bond geometry as shown.



RN 221247-54-7 HCAPLUS
CN .beta.-D-Glucopyranoside, hexadecyl 2-deoxy-2-(1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)-, 3,4,6-triacetate (9CI) (CA INDEX NAME)

Absolute stereochemistry.



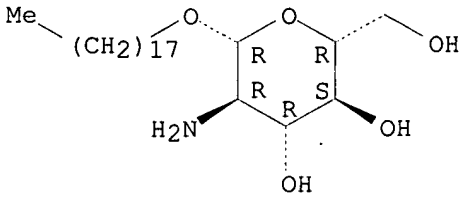
L13 ANSWER 1 OF 5 HCAPLUS COPYRIGHT 2001 ACS
ACCESSION NUMBER: 1999:590619 HCAPLUS
DOCUMENT NUMBER: 131:333590
TITLE: The folding and enzymatic activity of glucose oxidase in the glycolipid matrixes of different charges
AUTHOR(S): Li, J.-r.; Du, Y.-k.; Boullanger, P.; Jiang, L.
CORPORATE SOURCE: Institute of Photographic Chemistry, Molecular Science
Center, Laboratory of Colloid and Interface, Academia Sinica, Beijing, Peop. Rep. China
SOURCE: Thin Solid Films (1999), 352(1,2), 213-217
CODEN: THSFAP; ISSN: 0040-6090
PUBLISHER: Elsevier Science S.A.
DOCUMENT TYPE: Journal
LANGUAGE: English

AB The influence of lipid matrixes with different charges on the protein folding behavior has been investigated by using the Langmuir-Blodgett (LB) technique. To understand the conformation change under different forces, glucose oxidase (GOD) from Aspergillus Niger was used as a protein model. Eight glycolipids (1,2-o-dialkyl-3-o-.beta.-d-glycosylglycerols and alkyl 2-amino-2-deoxy-.beta.-d-glucopyranoside) were used as the matrixes for this investigation. It was obsd. that the GOD can penetrate into neutral glycolipid monolayer and change its conformation in favor of the .alpha.-helix formation. Moreover, GOD strongly adsorbed to the pos. charged glycolipid monolayer and change its conformation in favor of the .beta.-sheet formation. Enzymic activity measurements showed that the more the .alpha.-helix conformation content is in the GOD, the higher activity the GOD will be. This fact suggested a new way to mediate the conformation of protein in organized mol. assemblies, and provided a new thinking for the prepn. of biomimetic film and biosensor.

CC 6-3 (General Biochemistry)
Section cross-reference(s): 7
IT 81281-23-4 86363-39-5 86363-40-8 133128-66-2 159302-84-8 159302-85-9 **159405-31-9**
RL: BUU (Biological use, unclassified); PRP (Properties); BIOL (Biological study); USES (Uses)
(folding and enzymic activity of glucose oxidase in glycolipid matrixes

of different charges)
IT **159405-31-9**
RL: BUU (Biological use, unclassified); PRP (Properties); BIOL (Biological study); USES (Uses)
(folding and enzymic activity of glucose oxidase in glycolipid matrixes of different charges)
RN 159405-31-9 HCAPLUS
CN .beta.-D-Glucopyranoside, octadecyl 2-amino-2-deoxy- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

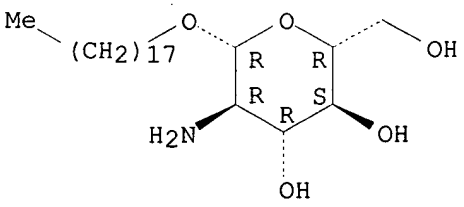


REFERENCE COUNT: 18
REFERENCE(S):
(2) Adler, A; Methods in Enzymology 1973, V27, P675 HCAPLUS
(6) Cheesman, D; Adv Protein Chem 1954, V9, P439 HCAPLUS
(7) Du, Y; Coll Surf B: Biointerfaces 1996, V7, P129 HCAPLUS
(9) Fenderson, B; BioEssays 1990, V12, P173 HCAPLUS
(10) Hakomori, S; Annu Rev Biochem 1981, V50, P733 HCAPLUS
ALL CITATIONS AVAILABLE IN THE RE FORMAT

L13 ANSWER 2 OF 5 HCAPLUS COPYRIGHT 2001 ACS
ACCESSION NUMBER: 1998:192924 HCAPLUS
DOCUMENT NUMBER: 128:305665
TITLE: Spectroscopic study of mixed Langmuir-Blodgett films of alkyl glycoside with positive charge and glucose oxidase
AUTHOR(S): Du, Yukou; Tang, Ji'an; Jiang, Long; Boullanger, Paul
CORPORATE SOURCE: Institute of Photographic Chemistry, Chinese Academy of Sciences, Beijing, 100101, Peop. Rep. China
SOURCE: Ganguang Kexue Yu Guang Huaxue (1998), 16(1), 32-37
CODEN: GKKHE9; ISSN: 1000-3231
PUBLISHER: Kexue Chubanshe
DOCUMENT TYPE: Journal
LANGUAGE: Chinese
AB Octadecyl-2-amino-2-deoxy-.beta.-D-glucopyranoside (C18) was synthesized. The interaction between C18 and glucose oxidase was studied by recording isotherms of monolayers. Mixed C18/glucose oxidase on Langmuir-Blodgett (LB) films was investigated by spectroscopy methods. From the results of CD spectrum, glucose oxidase immobilized in C18 films partly changed their secondary structure. Low temp. fluorescence studied showed that part of the glucose oxidase mols. in the LB film maintained its native structure,

and part became partly denatured.
CC 7-7 (Enzymes)
IT **159405-31-9**
RL: BUU (Biological use, unclassified); BIOL (Biological study); USES (Uses)
(glucose oxidase interaction with; spectroscopic study of mixed Langmuir-Blodgett films of alkyl glycoside with pos. charge and glucose
oxidase)
IT **159405-31-9**
RL: BUU (Biological use, unclassified); BIOL (Biological study); USES (Uses)
(glucose oxidase interaction with; spectroscopic study of mixed Langmuir-Blodgett films of alkyl glycoside with pos. charge and glucose
oxidase)
RN 159405-31-9 HCAPLUS
CN .beta.-D-Glucopyranoside, octadecyl 2-amino-2-deoxy- (9CI) (CA INDEX NAME)

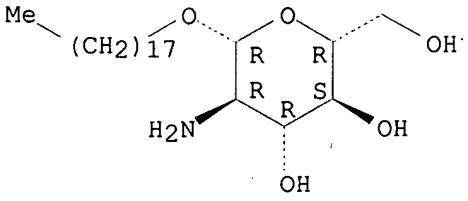
Absolute stereochemistry.



L13 ANSWER 3 OF 5 HCAPLUS COPYRIGHT 2001 ACS
ACCESSION NUMBER: 1995:977401 HCAPLUS
DOCUMENT NUMBER: 124:176728
TITLE: Synthesis and surface-active properties of some alkyl 2-amino-2-deoxy-.beta.-D-glucopyranosides
AUTHOR(S): Boullanger, Paul; Chevalier, Yves; Croizier, Marie-Christine; Lafont, Dominique; Sancho, Marie-Rose
CORPORATE SOURCE: Lab. Chimie Org., Univ. Lyon 1, Villeurbanne, F-69622, Fr.
SOURCE: Carbohydr. Res. (1995), 278(1), 91-101
CODEN: CRBRAT; ISSN: 0008-6215
DOCUMENT TYPE: Journal
LANGUAGE: English
AB Several alkyl 2-acetamido-2-deoxy-.beta.-D-glucopyranosides were synthesized using either the oxazoline or the N-allyloxycarbonyl procedure. The latter procedure gave better yields with fatty alcs. and cholesterol. The derivs. thus prepd. were partly or fully deprotected and
their surface-active properties assessed.
CC 33-7 (Carbohydrates)
Section cross-reference(s): 32, 46
IT 72205-17-5P 147025-06-7P 147126-58-7P 152914-68-6P 152914-69-7P 159302-81-5P 159302-82-6P 159302-85-9P **159405-31-9P**

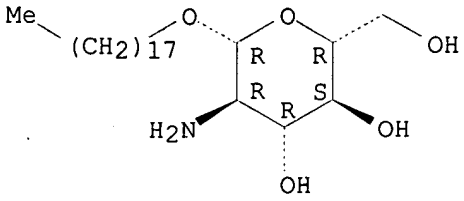
173725-28-5P 173725-29-6P 173934-01-5P
RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation)
(synthesis and surface-active properties of alkyl aminodeoxyglucopyranosides)
IT **159405-31-9P**
RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation)
(synthesis and surface-active properties of alkyl aminodeoxyglucopyranosides)
RN 159405-31-9 HCAPLUS
CN .beta.-D-Glucopyranoside, octadecyl 2-amino-2-deoxy- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



L13 ANSWER 4 OF 5 HCAPLUS COPYRIGHT 2001 ACS
ACCESSION NUMBER: 1994:712595 HCAPLUS
DOCUMENT NUMBER: 121:312595
TITLE: Mesomorphic amino sugars
AUTHOR(S): Stangier, P.; Vill, V.; Rohde, S.; Jeschke, U.; Thiem, J.
CORPORATE SOURCE: Dep. Chem., Univ. Alberta, Edmonton, AB, Can.
SOURCE: Liq. Cryst. (1994), 17(4), 589-95
CODEN: LICRE6; ISSN: 0267-8292
DOCUMENT TYPE: Journal
LANGUAGE: English
AB New liq. cryst. compds. were prepd. from glucosamine and 6-amino-6-deoxyhexopyranose. The monoalkylated carbohydrates show smectic phases. The influence of the amino group on the clearing temps. is minor.
The salts of the cyclic amines can form smectic or discotic mesophases. The clearing points are lower than those obsd. for acyclic amines.
CC 75-11 (Crystallography and Liquid Crystals)
Section cross-reference(s): 33
IT 159302-74-6P 159302-75-7P 159302-82-6P 159302-86-0P 159302-88-2P
159302-89-3P 159302-90-6P 159302-91-7P 159302-92-8P
159405-31-9P
RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation)
(prepn. and transition temps. of)
IT **159405-31-9P**
RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation)
(prepn. and transition temps. of)
RN 159405-31-9 HCAPLUS
CN .beta.-D-Glucopyranoside, octadecyl 2-amino-2-deoxy- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

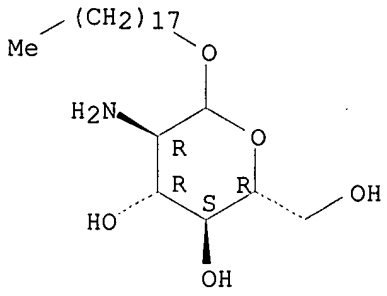


L13 ANSWER 5 OF 5 HCAPLUS COPYRIGHT 2001 ACS
ACCESSION NUMBER: 1991:415307 HCAPLUS
DOCUMENT NUMBER: 115:15307
TITLE: Skin-lightening cosmetics containing glucosamines
INVENTOR(S): Mishima, Yutaka; Okajima, Takehiko; Hori, Toshiro; Nishimoto, Katsuya; Oyama, Yasuaki
PATENT ASSIGNEE(S): Taiyo Chemical Co., Ltd., Japan
SOURCE: Jpn. Kokai Tokkyo Koho, 6 pp.
CODEN: JKXXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 02273608	A2	19901108	JP 1989-96859	19890417

OTHER SOURCE(S): MARPAT 115:15307
AB Skin-lightening cosmetics contain .gtoreq.1 glucosamines I (R1-5 = H, C<30 acyl, alkyl, alkenyl, alkynyl, aryl; .gtoreq.1 of R1-5 = alkyl, alkenyl, alkynyl, aryl) and/or their salts as active ingredients. The cosmetics are safe and prevent skin darkening caused by the sun light. A lotion comprised poly(oxyethylene)(20) monooleate 1.0, EtOH 3.0, polyethylene glycol-600 5.0, citric acid 0.03, Na citrate 0.2, 1-O-ethyltetraacetylglucosamine 0.1, methylparaben 0.1, fragrances, and H2O to 100 wt. parts.
IC ICM A61K007-00
CC 62-4 (Essential Oils and Cosmetics)
IT 134120-77-7 134227-31-9 **134227-32-0** 134227-33-1
134227-34-2 134275-62-0 134309-19-6
RL: BIOL (Biological study)
(skin-lightening cosmetics contg.)
IT **134227-32-0**
RL: BIOL (Biological study)
(skin-lightening cosmetics contg.)
RN 134227-32-0 HCAPLUS
CN D-Glucopyranoside, octadecyl 2-amino-2-deoxy-, hydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.



● HCl